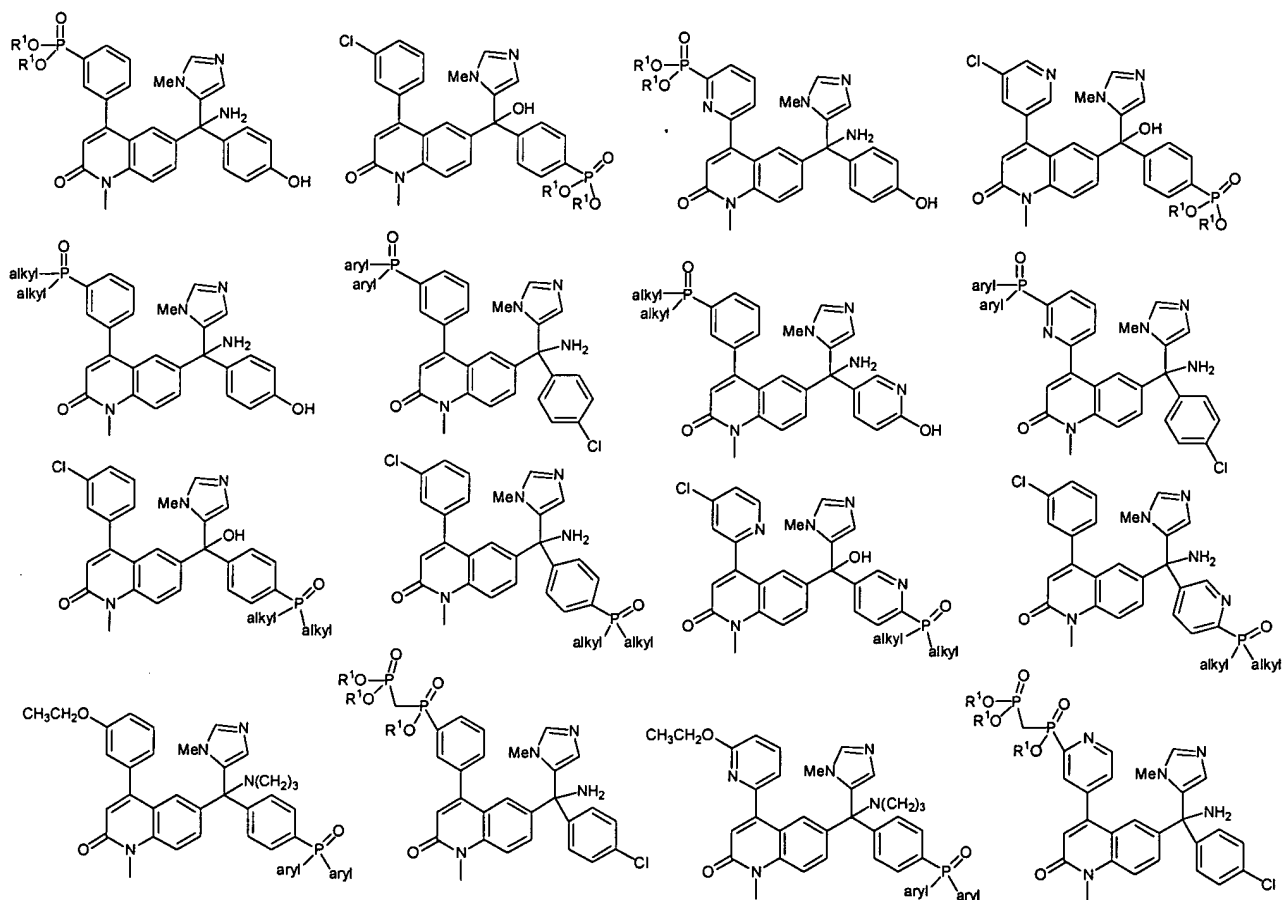




1. (Currently Amended) A compound chosen from the following:

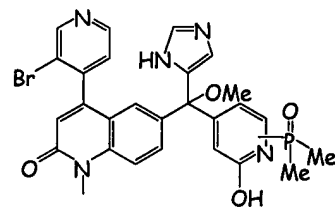
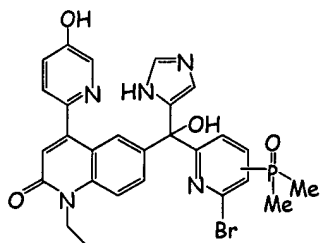
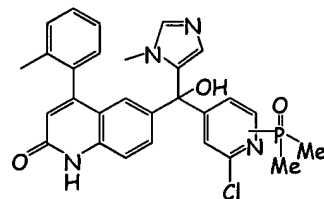
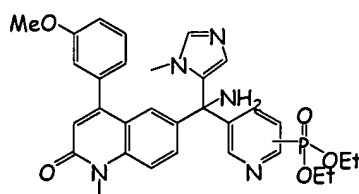
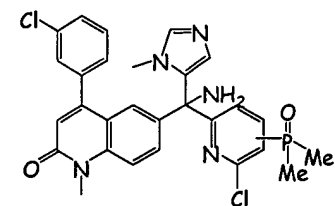
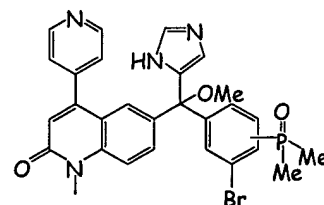
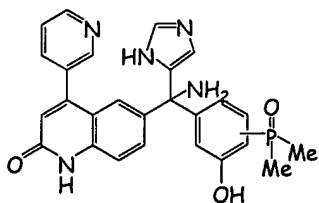
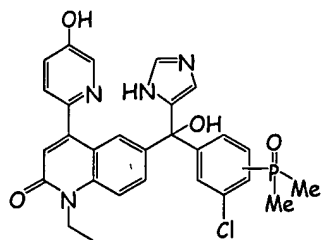
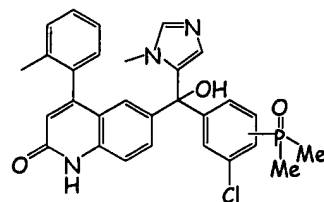
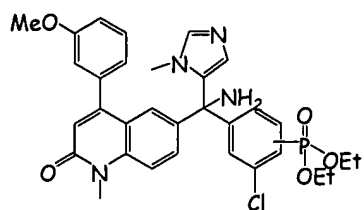
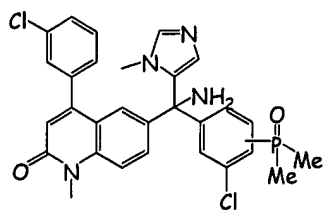


~~each occurrence of Y is independently O, S, or NR⁴, or a chemical bond linking R⁴ to P;~~

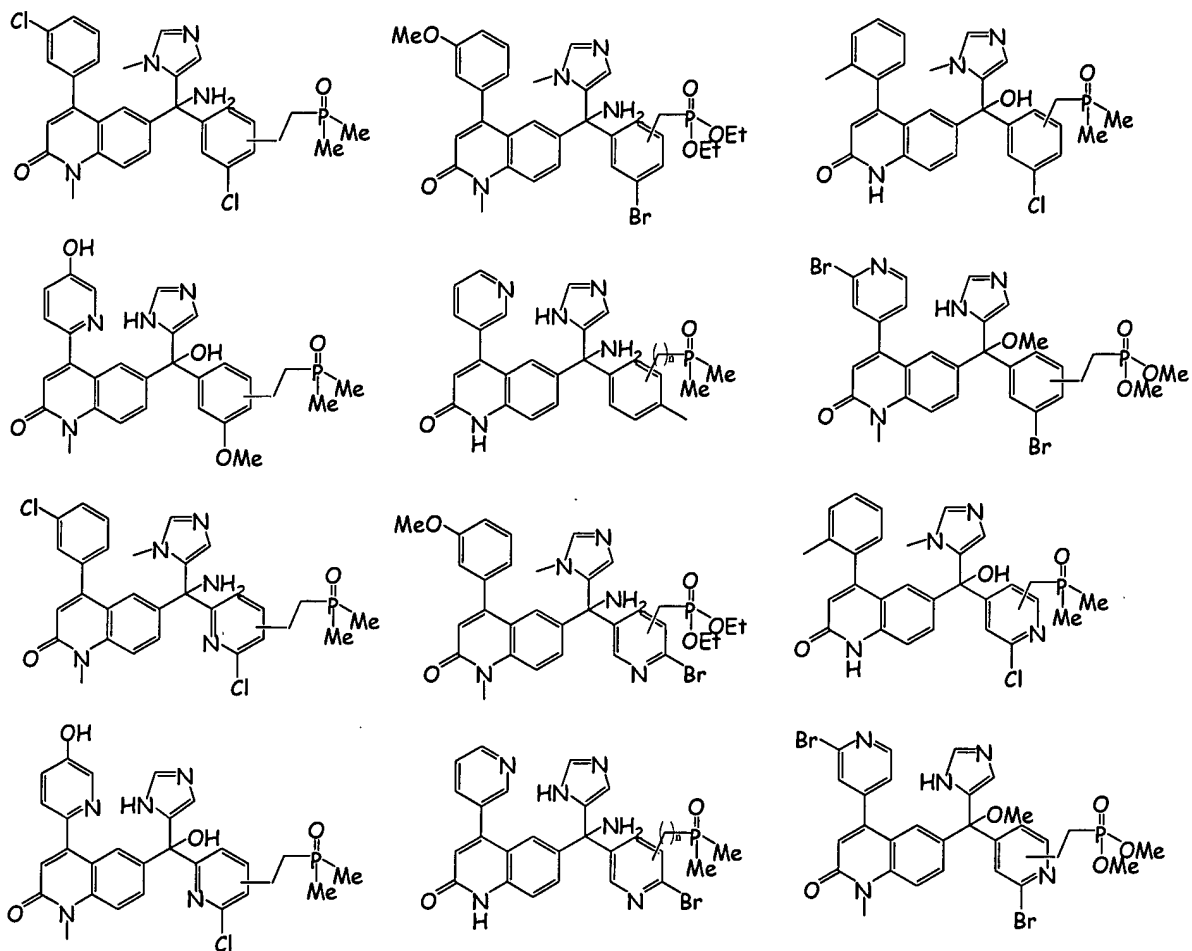
each occurrence of R¹ is independently a substituted or unsubstituted aliphatic, heteroaliphatic, aryl or heteroaryl moiety, or, except in YR¹ moieties in which Y is a covalent bond, R¹ may also be H; and

each alkyl and aryl group may be substituted or unsubstituted.

2. (Original) A compound chosen from the following:

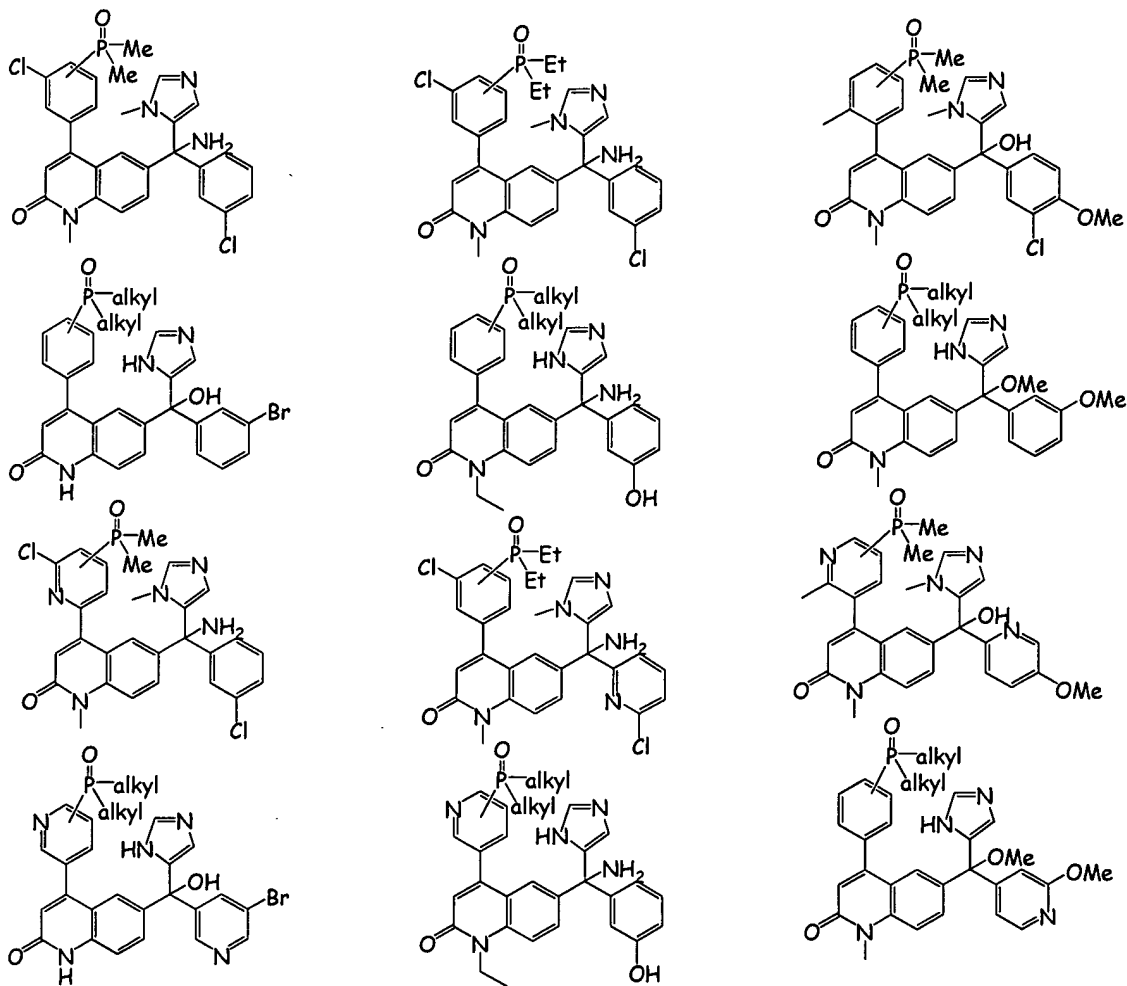


3. (Original) A compound chosen from:



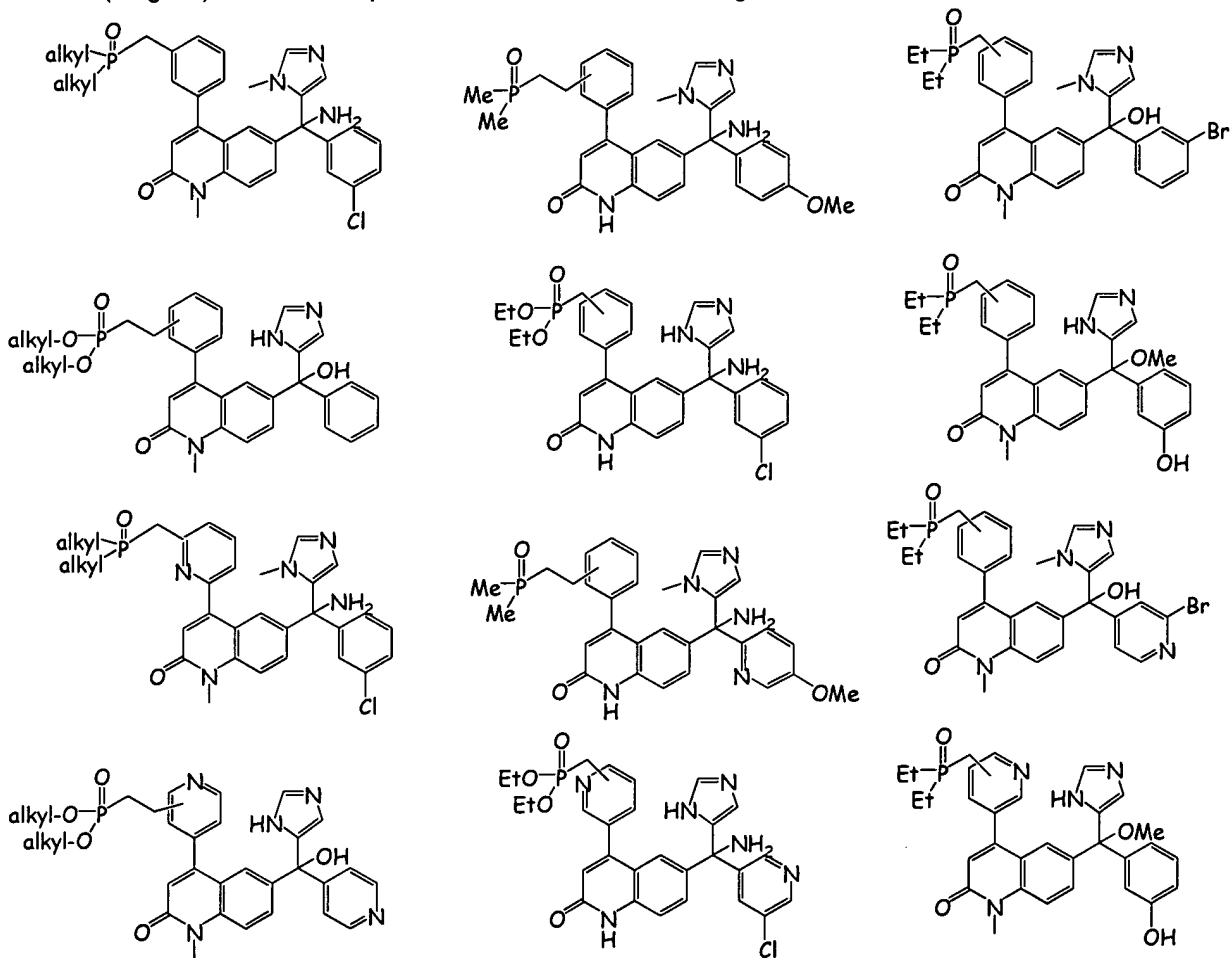
wherein n is an integer from 0 to 6.

4. (Original) A compound chosen from the following:



wherein each alkyl group may be substituted or unsubstituted.

5. (Original) A compound chosen from the following:



wherein each alkyl group may be substituted or unsubstituted.

6. (Original) The compound of any of claims 1 – 5 in which each alkyl group is a substituted or unsubstituted methyl, ethyl, n-propyl, isopropyl, cyclopropyl, -CH₂-cyclopropyl, n-butyl, sec-butyl, isobutyl, tert-butyl, cyclobutyl, -CH₂-cyclobutyl, n-pentyl, sec-pentyl, isopentyl, tert-pentyl, cyclopentyl, -CH₂-cyclopentyl, n-hexyl, sec-hexyl, cyclohexyl, or -CH₂-cyclohexyl moiety; or a benzyl or phenethyl moiety or a heteroaromatic or substituted derivative thereof.